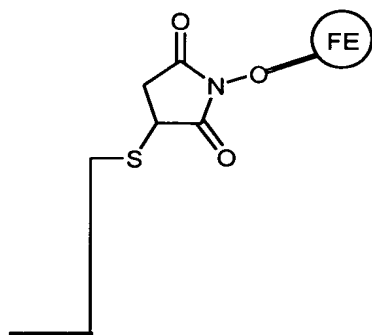


Amendments to the Claims:

This listing of claims will replace all prior versions, and listings, of claims in the application:

Listing of Claims:

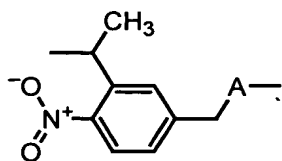
1. (Original) A building block of the general formula



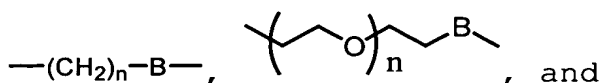
capable of transferring a functional entity (FE) to a recipient reactive group, wherein

the lower horizontal line is a **Complementing Element** identifying the functional entity and the vertical line between the complementing element and the S atom is a **Spacer**.

2. (Original) The building block of claim 1, wherein the spacer is a valence bond, C<sub>1</sub>-C<sub>6</sub> alkylene-A-, C<sub>1</sub>-C<sub>6</sub> alkenylene-A-, C<sub>2</sub>-C<sub>6</sub> alkynylene-A-, or

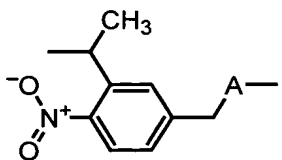


said spacer optionally being connected through A to a moiety selected from

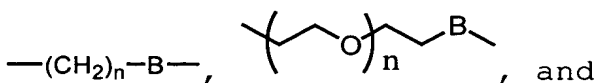


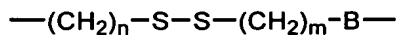
where A is a valence bond,  $-C(O)NR^1-$ ,  $-NR^1-$ ,  $-O-$ ,  $-S-$ , or  $-C(O)-O-$ ; B is a valence bond,  $-O-$ ,  $-S-$ ,  $-NR^1-$  or  $-C(O)NR^1-$  and connects to the S atom of the carrier;  $R^1$  is selected independently from H,  $C_1-C_6$  alkyl,  $C_3-C_7$  cycloalkyl,  $C_1-C_6$  alkylen-aryl, or aryl substituted with 0-5 halogen atoms selected from -F, -Cl, -Br and -I; and n and m independently are integers ranging from 1 to 10.

3. (Original) The compound according to claim 1, wherein the **Spacer** is  $C_1-C_6$  alkylen-A-,  $C_1-C_6$  alkenylene-A-,  $C_2-C_6$  alkynylene-A-, or



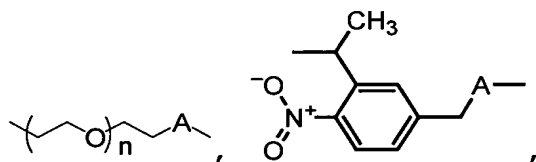
said spacer optionally being connected through A to a moiety selected from



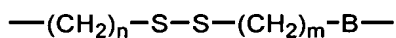


where A is  $-C(O)NR^1-$ , or  $-S-$ ; B is  $-S-$ ,  $-NR^1-$  or  $-C(O)NR^1-$  and connects to S-C-connecting group;  $R^1$  is selected independently from H,  $C_1-C_6$  alkyl,  $C_1-C_6$  alkylene-aryl, or aryl; and n and m independently are integers ranging from 1 to 6.

4. (Original) The compound according to claim 1, wherein **Spacer** is  $-A-$ , a group  $C_1-C_6$  alkylene- $A-$ ,  $C_2-C_6$  alkenylene- $A-$ , or  $C_2-C_6$  alkynylene- $A-$  optionally substituted with 1 to 3 hydroxy groups, or



said spacer being connected through A to a linker selected from

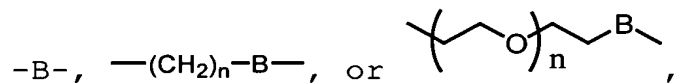


where A is a valence bond,  $-NR^2-$ ,  $-C(O)NR^2-$ ,  $-NR^2-C(O)-$ ,  $-O-$ ,  $-S-$ ,  $-C(O)-O-$  or  $-OP(=O)(O^-)-O-$ ; B is a valence bond,  $-O-$ ,  $-S-$ ,  $-NR^2-$ ,  $-C(O)-$  or  $-C(O)NR^2-$  and connects to S-C-connecting group;  $R^2$  is selected independently from H,  $C_1-C_6$  alkyl,  $C_3-C_7$

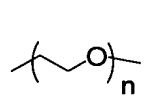
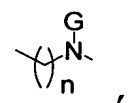
cycloalkyl, aryl,  $C_1-C_6$  alkylene-aryl,  $(CH_2CH_2O)_n$  or  $(CH_2CH_2N^G)_n$ ; G is

H or C<sub>1</sub>-C<sub>6</sub> alkyl; and n and m independently are integers ranging from 1 to 10.

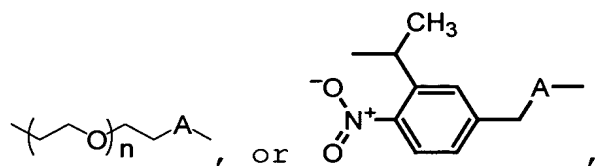
5. (Original) A compound according to claim 4, wherein the **spacer** is C<sub>2</sub>-C<sub>6</sub> alkenylene-A, said spacer being connected through A to a moiety selected from



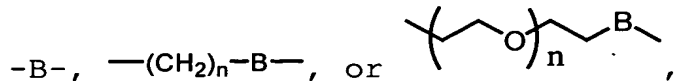
where A is a valence bond, -C(O)NR<sup>2</sup>-, -NR<sup>2</sup>-C(O)-, -S-, -C(O)-O- or -OP(=O)(O<sup>-</sup>)-O-; B is a valence bond, -S-, -NR<sup>2</sup>-, or -C(O)- and connects to S-C-connecting group; n and m independently are integers ranging from 1 to 10 and

R<sup>2</sup> is selected independently from H,  or , wherein G is H or C<sub>1</sub>-C<sub>6</sub> alkyl; and the spacer is connected to the complementing element through a nucleobase.

6. (Original) A compound according to claim 4, wherein the **spacer** is -A-,

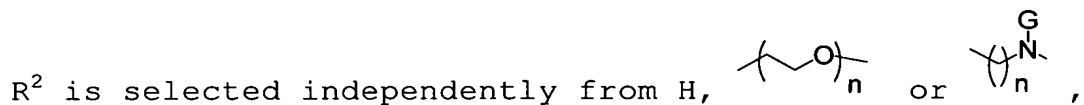


said spacer being connected through A to a moiety selected from



where A is a valence bond,  $-\text{NR}^2\text{-C(O)-}$ ,  $-\text{O-}$ , or  $-\text{S-}$ ; B is a valence bond,  $-\text{S-}$ ,  $-\text{NR}^2\text{-}$ , or  $-\text{C(O)-}$  and connects to S-C-connecting group;

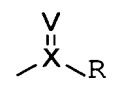
n and m independently are integers ranging from 1 to 10 and



wherein G is H or  $\text{C}_1\text{-C}_6$  alkyl; and the spacer is connected to the complementing element via a phosphorus group.

7. (Original) A compound according to claim 6, wherein the phosphorus group is a phosphate or thiophosphate group attached to a 3' or 5' end of a complementing element.

8. (Currently Amended) The building block according to ~~any of~~

~~the claims 1 to 7~~ claim 1, wherein **FE** is  where

X =  $-\text{C-}$ ,  $-\text{S-}$ ,  $-\text{P-}$ ,  $-\text{S(O)-}$ , or  $-\text{P(O)-}$ ,

V = O, S, NH, or  $\text{N-C}_1\text{-C}_6$  alkyl, and

R is H or selected among the group consisting of a  $\text{C}_1\text{-C}_6$  alkyl,  $\text{C}_2\text{-C}_6$  alkenyl,  $\text{C}_2\text{-C}_6$  alkynyl,  $\text{C}_4\text{-C}_8$  alkadienyl,  $\text{C}_3\text{-C}_7$  cycloalkyl,  $\text{C}_3\text{-C}_7$  cycloheteroalkyl, aryl, and heteroaryl, said group being substituted with 0-3  $\text{R}^4$ , 0-3  $\text{R}^5$  and 0-3  $\text{R}^9$  or  $\text{C}_1\text{-C}_3$  alkylene- $\text{NR}^4_2$ ,  $\text{C}_1\text{-C}_3$  alkylene- $\text{NR}^4\text{C(O)R}^8$ ,  $\text{C}_1\text{-C}_3$

alkylene-NR<sup>4</sup>C(O)OR<sup>8</sup>, C<sub>1</sub>-C<sub>2</sub> alkylene-O-NR<sup>4</sup><sub>2</sub>, C<sub>1</sub>-C<sub>2</sub>  
alkylene-O-NR<sup>4</sup>C(O)R<sup>8</sup>, C<sub>1</sub>-C<sub>2</sub> alkylene-O-NR<sup>4</sup>C(O)OR<sup>8</sup> substituted  
with 0-3 R<sup>9</sup>[[.]]<sub>1</sub>

where R<sup>4</sup> is H or selected independently among the  
group consisting of C<sub>1</sub>-C<sub>6</sub> alkyl, C<sub>2</sub>-C<sub>6</sub> alkenyl, C<sub>2</sub>-C<sub>6</sub> alkynyl,  
C<sub>3</sub>-C<sub>7</sub> cycloalkyl, C<sub>3</sub>-C<sub>7</sub> cycloheteroalkyl, aryl, heteroaryl, said  
group being substituted with 0-3 R<sup>9</sup> and

R<sup>5</sup> is selected independently from -N<sub>3</sub>, -CNO,  
-C(NOH)NH<sub>2</sub>, -NHOH, -NHNHR<sup>6</sup>, -C(O)R<sup>6</sup>, -SnR<sup>6</sup><sub>3</sub>, -B(OR<sup>6</sup>)<sub>2</sub>,  
-P(O)(OR<sup>6</sup>)<sub>2</sub> or the group consisting of C<sub>2</sub>-C<sub>6</sub> alkenyl, C<sub>2</sub>-C<sub>6</sub>  
alkynyl, C<sub>4</sub>-C<sub>8</sub> alkadienyl said group being substituted with 0-2  
R<sup>7</sup>,

where R<sup>6</sup> is selected independently from H, C<sub>1</sub>-C<sub>6</sub>  
alkyl, C<sub>3</sub>-C<sub>7</sub> cycloalkyl, aryl or C<sub>1</sub>-C<sub>6</sub> alkylene-aryl substituted  
with 0-5 halogen atoms selected from -F, -Cl, -Br, and -I; and  
R<sup>7</sup> is independently selected from -NO<sub>2</sub>, -COOR<sup>6</sup>, -COR<sup>6</sup>, -CN,  
-OSiR<sup>6</sup><sub>3</sub>, -OR<sup>6</sup> and -NR<sup>6</sup><sub>2</sub>[[.]]<sub>1</sub>

R<sup>8</sup> is H, C<sub>1</sub>-C<sub>6</sub> alkyl, C<sub>2</sub>-C<sub>6</sub> alkenyl, C<sub>2</sub>-C<sub>6</sub> alkynyl, C<sub>3</sub>-C<sub>7</sub>  
cycloalkyl, aryl or C<sub>1</sub>-C<sub>6</sub> alkylene-aryl substituted with 0-3  
substituents independently selected from -F, -Cl, -NO<sub>2</sub>, -R<sup>3</sup>,  
-OR<sup>3</sup>, -SiR<sup>3</sup><sub>3</sub>

R<sup>9</sup> is =O, -F, -Cl, -Br, -I, -CN, -NO<sub>2</sub>, -OR<sup>6</sup>, -NR<sup>6</sup><sub>2</sub>, -NR<sup>6</sup>-C(O)R<sup>8</sup>,  
-NR<sup>6</sup>-C(O)OR<sup>8</sup>, -SR<sup>6</sup>, -S(O)R<sup>6</sup>, -S(O)<sub>2</sub>R<sup>6</sup>, -COOR<sup>6</sup>, -C(O)NR<sup>6</sup><sub>2</sub> and  
-S(O)<sub>2</sub>NR<sup>6</sup><sub>2</sub>.

9. (Original) A compound according to claim 8, wherein R is H or selected among the group consisting of a C<sub>1</sub>-C<sub>6</sub> alkyl, C<sub>2</sub>-C<sub>6</sub> alkenyl, C<sub>2</sub>-C<sub>6</sub> alkynyl, C<sub>4</sub>-C<sub>8</sub> alkadienyl, C<sub>3</sub>-C<sub>7</sub> cycloalkyl, C<sub>3</sub>-C<sub>7</sub> cycloheteroalkyl, aryl, and heteroaryl, said group being substituted with 0-3 R<sup>5</sup> and 0-3 R<sup>9</sup>, or selected among the group consisting of C<sub>1</sub>-C<sub>3</sub> alkylene-NR<sup>4</sup><sub>2</sub>, C<sub>1</sub>-C<sub>3</sub> alkylene-NR<sup>4</sup>C(O)R<sup>8</sup>, C<sub>1</sub>-C<sub>3</sub> alkylene-NR<sup>4</sup>C(O)OR<sup>8</sup>, C<sub>1</sub>-C<sub>2</sub> alkylene-O-NR<sup>4</sup><sub>2</sub>, C<sub>1</sub>-C<sub>2</sub> alkylene-O-NR<sup>4</sup>C(O)R<sup>8</sup>, and C<sub>1</sub>-C<sub>2</sub> alkylene-O-NR<sup>4</sup>C(O)OR<sup>8</sup> substituted with 0-3 R<sup>9</sup>.

10. (Currently Amended) A compound according to ~~claims 8 or 9~~ claim 8, wherein R is H or selected among the group consisting of C<sub>1</sub>-C<sub>6</sub> alkyl, C<sub>2</sub>-C<sub>6</sub> alkenyl, C<sub>2</sub>-C<sub>6</sub> alkynyl, C<sub>4</sub>-C<sub>8</sub> alkadienyl, C<sub>3</sub>-C<sub>7</sub> cycloalkyl, C<sub>3</sub>-C<sub>7</sub> cycloheteroalkyl, aryl, and heteroaryl, said group being substituted with 0-3 R<sup>5</sup> and 0-3 R<sup>9</sup>.

11. (Currently Amended) A compound according to ~~any of the claims 8 to 10~~ claim 8, wherein R is selected among the group consisting of C<sub>1</sub>-C<sub>3</sub> alkylene-NR<sup>4</sup><sub>2</sub>, C<sub>1</sub>-C<sub>3</sub> alkylene-NR<sup>4</sup>C(O)R<sup>8</sup>, C<sub>1</sub>-C<sub>3</sub> alkylene-NR<sup>4</sup>C(O)OR<sup>8</sup>, C<sub>1</sub>-C<sub>2</sub> alkylene-O-NR<sup>4</sup><sub>2</sub>, C<sub>1</sub>-C<sub>2</sub> alkylene-O-NR<sup>4</sup>C(O)R<sup>8</sup>, and C<sub>1</sub>-C<sub>2</sub> alkylene-O-NR<sup>4</sup>C(O)OR<sup>8</sup> substituted with 0-3 R<sup>9</sup>.

12. (Currently Amended) A compound according to ~~any of the claims 1 to 11~~ claim 1, wherein X = C and V = O or S.

13. (Currently Amended) A compound according to ~~claims 1 to 12~~ claim 1, wherein X = C and V = O.

14. (Currently Amended) A compound according to ~~claims 1 to 13~~ claim 1, wherein complementing element is a nucleic acid.

15. (Currently Amended) A compound according to ~~claims 1 to 14~~ claim 1, wherein Complementing element is a sequence of nucleotides selected from the group of DNA, RNA, LNA PNA, or morpholino derivatives.

16. (Currently Amended) A library of compounds according to ~~any of the claims 1 to 15~~ claim 1, wherein each different member of the library comprises a complementing element having a unique sequence of nucleotides, which identifies the functional entity.

17. (Currently Amended) A method for transferring a functional entity to a recipient reactive group, comprising the steps of



providing one or more building blocks according to  
~~claims 1 to 15~~ claim 1,

contacting the one or more building blocks with a  
corresponding encoding element associated with a recipient  
reactive group under conditions which allow for a recognition  
between the one or more complementing elements and the coding  
elements, said contacting being performed prior to,  
simultaneously with, or subsequent to a transfer of the  
functional entity to the recipient reactive group.

18. (Original) The method according to claim 17, wherein the  
coding element comprises one or more coding sequences  
comprised of 1 to 50 nucleotides and the one or more  
complementing elements comprises a sequence of nucleotides  
complementary to one or more of the coding sequences.

19. (Currently Amended) The method of ~~claims 17 or 18~~ claim  
17, wherein the recipient reactive group is an amine group,  
which may be part of a chemical scaffold, and the linkage  
between the functional entity and the scaffold is of the  
general chemical structure:

**Scaffold-NH-X(=V) -R**

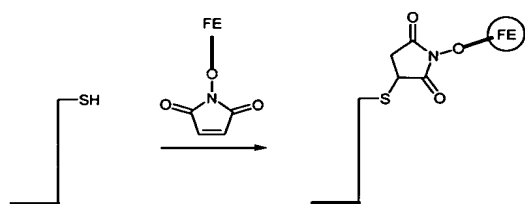
In which

X = -C-, -S-, -P-, -S(O)-, -P(O)-, and

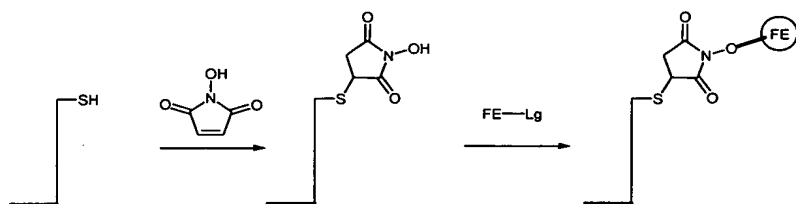
V = O, S, NH, N-C<sub>1</sub>-C<sub>6</sub> alkyl.

20. (Original) The method according to claim 19, wherein X is C and V is O.

21. (Original) A process for preparing a building block according to claim 1, comprising the step of

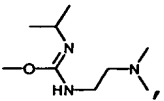
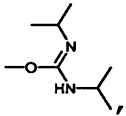
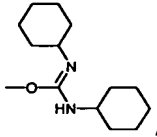


22. (Original) A process for preparing a building block according to claim 1, comprising the steps of



where Lg is a leaving group.

23. (Original) A process according to claim 18, wherein the

leaving group is selected from , , , Cl,  
Br.